Journal of The Chemical Society

B

Physical Organic

Indexes

1966

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LONDON: THE CHEMICAL SOCIETY
1966

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Richard Clay (The Chaucer Press), Ltd.,
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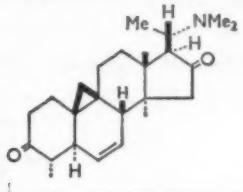
ERRATA prior to 1 January 1966

Vol. 1948, page 239, line 10*. *For* m. p. 223—224°, $[\alpha]_D^{18}$ −16·4° *read* m. p. 223—224°, $[\alpha]_D^{18}$ +16·4°.

Vol. 1965, page 4513, Table, line 3*. *For* *N*-methyl groups at C-3 and C-20 *read* *N*-methyl groups at C-20 and C-3.

Vol. 1965, page 4521. Formula (XXX) should be shown as

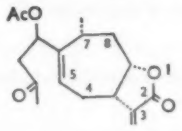


Vol. 1965, page 6548, line 23. *For* 72% of 4-hydroxycyclohexyl phosphate and, *read* 72% of cyclohexane diphosphate and

---

Vol. 1965, page 6777, Paper 1255, Title line 2. *For* Synthesis of Poly-β-L- and, *read* Synthesis of Poly-γ-L- and.

Vol. 1965, page 7009. Formula (II) should be shown as



Vol. 1965, page 7174, Table 5, compound 36. *For* $C_{13}H_{18}ClN_3$ *read* $C_{13}H_{19}ClN_2$.

Vol. 1965, page 7180, Table footnote* line 3. (*a*) A neutral by-product was indol-3-ylacetaldoxime (VI), *delete* (VI).

Vol. 1965, page 7425, Summary line 4. *For* a 3,4-cyclic product *read* a 2,4-cyclic product.

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* From foot of main text.

ERRATA. Section B

1966, page 56, Summary line 2. *For* pH 1·5—9·4 *read* 1·5—9·3.

1966, page 57, Table 1. The figure of 1·85 in the column headed pK_a' should be lowered one line.

1966, page 59, Table 3. Under N-*Cyclohexylhydroxylamine with n-butyraldehyde*. *For* A_{a1} *read* K_{a1}. *For* A_h *read* K_h.
Under N-*Cyclohexylhydroxylamine with pivalaldehyde*. *For* K_{Q1} *read* K_{a1}. *For* $K_h = 0$ *read* $K_h \sim 0$.

1966, page 60, Table 5, column headed k_{H^+}, line 10. *For* $6{\cdot}24 \times 10^4$ *read* $6{\cdot}29 \times 10^4$.

1966, page 61. Equation 13. *For* $K\begin{matrix}\overset{(II)}{\rightleftharpoons}\\\underset{(III)}{}\end{matrix}H^+$ *read* $K\begin{matrix}\overset{(II)}{\rightleftharpoons}\\\underset{(III)}{}\end{matrix}-H^+$

1966, page 76, Table 2, column headed CH_3. Compound (XI) *delete* 6·11(s). Compound (XII) *insert* 6·11(s) after the figure 8·89(s). Compound (XV) *For* 8·89(s) *read* 8·39(s).

1966, page 171, Table 3. The authors state " The result is that the K_c values in table 3 should be *half* the values quoted, and the ε values *twice* the figures given ".

1966, page 209, Summary line 2 and Results line 2. *For trans*-hex-4-enal. *read* 4-methylpent-4-enal.

1966, page 210, right hand column, line 6. *For* in the present case *read* from the 1-methylallyl compound.[6] *Insert* ref. 6 P. Cresson, *Bull. Soc. chim. France*, 1964, 2629.

1966, page 503, left hand column, line 19*. *For* of 2,4-dinitroanisole [10] *read* of 2,4-dinitroanisole [2].

1966, page 503, right hand column, line 13*. *For* position than nitrobenzene *read* position than trinitrobenzene.

1966, page 557, right hand column, line 30. *For* Tropane (6·5 g.) *read* Tropanone (6·5 g.).

1966, page 559, Equation (9) should read

$$0{\cdot}79k_a/(0{\cdot}20k_e + 0{\cdot}01k_a) = 2 \text{ or } 0{\cdot}5 \qquad (9)$$

1966, page 560, line 6. Rate ($\times 10^4$) figures should read as follows: 1·14 13·5 80·7 397·9

1966, page 949 and 954. *Interchange figures as set out*
" The mass spectrum of 3-methoxythioanisole reproduced as Figure 4 in Part XIII, should appear as Figure 9 in Part XIV. Conversely, the mass spectrum of diphenyl disulphide, reproduced as Figure 9 in Part XIV, should appear as Figure 4 in Part XIII."

1966, page 1057, right hand column, line 3. After molal transition energy, E_T* *insert*

* A better correction exists with the refractive index of the solvent indicating that polarisability of the solvent is important in stabilising the excited state of the complex: (H. M. Rosenberg, *Chem. Comm.* 1965, 312: H. M. Rosenberg, E. Eimutis, and D. Hale, *Canad. J. Chem.*, 1966, 44, 2405.).

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* From foot of main text.

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Printed in Great Britain by Richard Clay (The Chaucer Press), Ltd.
Bungay, Suffolk